

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100324\
 Data File : BM047826.D
 Acq On : 04 Oct 2024 03:11
 Operator : RC/JU
 Sample : P4084-17
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DCYM1

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 10/04/2024
 Supervised By :mohammad ahmed 10/06/2024

Quant Time: Oct 04 04:00:14 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Sep 28 02:30:36 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.804	152	280264	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.592	136	1104900	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.439	164	676162	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.174	188	1351199	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.397	240	1295566	20.000	ng/u1	-0.01
88) Perylene-d12	24.397	264	1506053	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	26838	3.095	ng/uL	0.00
4) Pyridine-d5	3.681	84	27160	1.070	ng/u1	0.00
7) Phenol-d5	6.969	99	561242	21.178	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	364654	19.391	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.334	132	463377	24.032	ng/u1	-0.01
15) 4-Methylphenol-d8	8.504	113	459394	23.315	ng/u1	-0.01
21) Nitrobenzene-d5	8.957	128	238288	27.087	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.675	143	227664	27.005	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	10.216	165	445087	26.429	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.727	131	424424m	16.301	ng/u1	-0.01
46) Dimethylphthalate-d6	13.845	166	1219727	24.881	ng/u1	-0.01
49) Acenaphthylene-d8	14.127	160	1663969	28.628	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.627	143	200461	20.538	ng/u1	0.00
60) Fluorene-d10	15.427	176	1259386	29.659	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.539	200	144479	18.272	ng/u1	-0.01
73) Anthracene-d10	17.274	188	1961877	29.529	ng/u1	-0.01
81) Pyrene-d10	19.556	212	2287946	31.116	ng/u1	-0.01
92) Benzo(a)pyrene-d12	24.191	264	2377454	30.425	ng/u1	-0.02

Target Compounds Qvalue

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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