

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082223\
 Data File : BP016699.D
 Acq On : 22 Aug 2023 12:18
 Operator : MA/JU
 Sample : 04037-18
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZYN7

Quant Time: Aug 23 00:43:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Aug 17 01:38:34 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.046	152	320572	20.000	ng/u1	0.00
20) Naphthalene-d8	10.875	136	1461292	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.693	164	933926	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.463	188	2149655	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	1915574	20.000	ng/u1	0.00
88) Perylene-d12	24.139	264	1896527	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	38370	4.548	ng/uL	0.00
4) Pyridine-d5	3.828	84	136067	5.301	ng/u1	0.00
7) Phenol-d5	7.181	99	140563	4.756	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.381	67	436002	22.398	ng/u1	0.00
11) 2-Chlorophenol-d4	7.563	132	374509	17.643	ng/u1	0.00
15) 4-Methylphenol-d8	8.746	113	259084	11.025	ng/u1	-0.01
21) Nitrobenzene-d5	9.240	128	252459	24.553	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.957	143	252534	25.859	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.481	165	410159	21.342	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.028	131	611082	18.532	ng/u1	-0.01
46) Dimethylphthalate-d6	14.104	166	1587819	23.506	ng/u1	0.00
49) Acenaphthylene-d8	14.393	160	1809211	23.729	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.898	143	45946	3.339	ng/u1	0.00
60) Fluorene-d10	15.687	176	1404522	24.389	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.810	200	184034	17.996	ng/u1	0.00
73) Anthracene-d10	17.563	188	2398783	25.798	ng/u1	0.00
81) Pyrene-d10	19.792	212	2806268	27.959	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.968	264	2313544	25.455	ng/u1	0.00

Target Compounds	Qvalue

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

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