

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092523\
 Data File : BP017351.D
 Acq On : 26 Sep 2023 08:52
 Operator : MA/JU
 Sample : 04472-11
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EZYL0

Quant Time: Sep 27 00:25:36 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Sep 26 04:52:29 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	181472	20.000	ng/u1	0.00
20) Naphthalene-d8	10.746	136	770123	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.586	164	470362	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.357	188	1050769	20.000	ng/u1	0.00
79) Chrysene-d12	21.463	240	984025	20.000	ng/u1	0.00
88) Perylene-d12	23.974	264	1111647	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	19946	4.515	ng/uL	0.00
4) Pyridine-d5	3.728	84	327634	26.526	ng/u1	0.00
7) Phenol-d5	7.081	99	425096	28.522	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	260258	28.935	ng/u1	0.00
11) 2-Chlorophenol-d4	7.446	132	360752	30.518	ng/u1	0.00
15) 4-Methylphenol-d8	8.640	113	309441	26.705	ng/u1	0.00
21) Nitrobenzene-d5	9.122	128	175995	31.611	ng/u1	0.00
24) 2-Nitrophenol-d4	9.840	143	181968	30.101	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	370619	32.450	ng/u1	0.00
31) 4-Chloroaniline-d4	10.910	131	452512	27.351	ng/u1	0.00
46) Dimethylphthalate-d6	14.004	166	1244814	36.943	ng/u1	0.00
49) Acenaphthylene-d8	14.287	160	1398477	36.086	ng/u1	0.00
54) 4-Nitrophenol-d4	14.816	143	74548	12.699	ng/u1	0.00
60) Fluorene-d10	15.586	176	980572	33.803	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.716	200	32868	5.550	ng/u1	0.00
73) Anthracene-d10	17.457	188	1758013	37.724	ng/u1	0.00
81) Pyrene-d10	19.704	212	2157199	40.539	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.810	264	2038459	38.086	ng/u1	0.00

Target Compounds	Qvalue

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

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