

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082323\
 Data File : BP016713.D
 Acq On : 23 Aug 2023 14:08
 Operator : MA/JU
 Sample : PB155004BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK004

Quant Time: Aug 24 01:05:49 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Aug 23 18:42:31 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.040	152	422173	20.000	ng/u1	0.00
20) Naphthalene-d8	10.869	136	1877692	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.693	164	1145603	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.463	188	2198701	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	1344850	20.000	ng/u1	0.00
88) Perylene-d12	24.133	264	1515727	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	91973	8.278	ng/uL	0.00
4) Pyridine-d5	3.823	84	1102432	32.614	ng/u1	0.00
7) Phenol-d5	7.181	99	1265563	32.516	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.381	67	827063	32.263	ng/u1	0.00
11) 2-Chlorophenol-d4	7.558	132	907087	32.448	ng/u1	0.00
15) 4-Methylphenol-d8	8.746	113	978403	31.615	ng/u1	0.00
21) Nitrobenzene-d5	9.240	128	469104	35.506	ng/u1	0.00
24) 2-Nitrophenol-d4	9.957	143	494475	39.405	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.475	165	834332	33.786	ng/u1	0.00
31) 4-Chloroaniline-d4	11.028	131	1316643	31.074	ng/u1	0.00
46) Dimethylphthalate-d6	14.104	166	2759704	33.305	ng/u1	0.00
49) Acenaphthylene-d8	14.393	160	3137450	33.547	ng/u1	0.00
54) 4-Nitrophenol-d4	14.893	143	359859	21.319	ng/u1	0.00
60) Fluorene-d10	15.687	176	2294678	32.484	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.810	200	252928	24.181	ng/u1	0.00
73) Anthracene-d10	17.563	188	3256259	34.239	ng/u1	0.00
81) Pyrene-d10	19.792	212	3030592	43.007	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.963	264	2532356	34.862	ng/u1	0.00

Target Compounds Qvalue

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

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