

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082123\
 Data File : BP016669.D
 Acq On : 21 Aug 2023 11:39
 Operator : MA/JU
 Sample : PB154934BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK934

Quant Time: Aug 22 01:05:19 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	366255	20.000	ng/u1	0.00
20) Naphthalene-d8	10.875	136	1675730	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.698	164	1051157	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.463	188	2204518	20.000	ng/u1	0.00
79) Chrysene-d12	21.563	240	1280600	20.000	ng/u1	0.00
88) Perylene-d12	24.139	264	1305881	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	73585	7.634	ng/uL	0.00
4) Pyridine-d5	3.828	84	920975	31.406	ng/u1	0.00
7) Phenol-d5	7.187	99	1039086	30.773	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.381	67	667964	30.035	ng/u1	0.00
11) 2-Chlorophenol-d4	7.563	132	740384	30.528	ng/u1	0.00
15) 4-Methylphenol-d8	8.752	113	813364	30.294	ng/u1	0.00
21) Nitrobenzene-d5	9.246	128	385636	32.706	ng/u1	0.00
24) 2-Nitrophenol-d4	9.963	143	406753	36.321	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.487	165	715555	32.468	ng/u1	0.00
31) 4-Chloroaniline-d4	11.034	131	1141275	30.182	ng/u1	0.00
46) Dimethylphthalate-d6	14.110	166	2428762	31.945	ng/u1	0.00
49) Acenaphthylene-d8	14.398	160	2762140	32.187	ng/u1	0.00
54) 4-Nitrophenol-d4	14.898	143	352701	22.772	ng/u1	0.00
60) Fluorene-d10	15.692	176	2049893	31.626	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.810	200	225292	21.482	ng/u1	0.00
73) Anthracene-d10	17.563	188	3100969	32.520	ng/u1	0.00
81) Pyrene-d10	19.798	212	3048488	45.432	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.968	264	2073005	33.124	ng/u1	0.00

Target Compounds Qvalue

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

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