

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082223\
 Data File : BP016695.D
 Acq On : 22 Aug 2023 09:45
 Operator : MA/JU
 Sample : PB154972BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK972

Quant Time: Aug 23 00:51:44 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	334140	20.000	ng/u1	0.00
20) Naphthalene-d8	10.875	136	1525465	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.692	164	979056	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.463	188	2243165	20.000	ng/u1	0.00
79) Chrysene-d12	21.557	240	1905104	20.000	ng/u1	0.00
88) Perylene-d12	24.145	264	1866735	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.387	96	66655	7.580	ng/uL	0.00
4) Pyridine-d5	3.822	84	826475	30.892	ng/u1	0.00
7) Phenol-d5	7.181	99	932766	30.279	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.381	67	616007	30.361	ng/u1	0.00
11) 2-Chlorophenol-d4	7.563	132	665472	30.077	ng/u1	0.00
15) 4-Methylphenol-d8	8.746	113	729327	29.775	ng/u1	-0.01
21) Nitrobenzene-d5	9.240	128	347595	32.384	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.957	143	365833	35.885	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.481	165	634647	31.633	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.028	131	1052877	30.587	ng/u1	-0.01
46) Dimethylphthalate-d6	14.104	166	2325423	32.838	ng/u1	0.00
49) Acenaphthylene-d8	14.392	160	2575186	32.219	ng/u1	-0.01
54) 4-Nitrophenol-d4	14.892	143	395140	27.391	ng/u1	0.00
60) Fluorene-d10	15.687	176	1975422	32.722	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.810	200	273143	25.596	ng/u1	0.00
73) Anthracene-d10	17.563	188	3209922	33.082	ng/u1	0.00
81) Pyrene-d10	19.798	212	3643602	36.501	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.968	264	2982384	33.337	ng/u1	0.00

Target Compounds	Qvalue

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

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