

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082823\
 Data File : BP016786.D
 Acq On : 28 Aug 2023 10:23
 Operator : MA/JU
 Sample : PB155031BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK031

Quant Time: Aug 29 00:55:56 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082423.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 25 01:47:30 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.028	152	364348	20.000	ng/u1	-0.01
20) Naphthalene-d8	10.857	136	1620411	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.687	164	1056149	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.457	188	2321084	20.000	ng/u1	0.00
79) Chrysene-d12	21.551	240	1617875	20.000	ng/u1	0.00
88) Perylene-d12	24.127	264	1424434	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	70931	8.046	ng/uL	0.00
4) Pyridine-d5	3.811	84	839331	31.155	ng/u1	-0.01
7) Phenol-d5	7.169	99	976497	30.141	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.363	67	612681	29.672	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.546	132	749811	30.782	ng/u1	-0.01
15) 4-Methylphenol-d8	8.734	113	778026	29.418	ng/u1	-0.01
21) Nitrobenzene-d5	9.228	128	378655	30.906	ng/u1	-0.01
24) 2-Nitrophenol-d4	9.946	143	405174	31.374	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.469	165	730185	30.328	ng/u1	0.00
31) 4-Chloroaniline-d4	11.016	131	1069228	28.959	ng/u1	-0.01
46) Dimethylphthalate-d6	14.098	166	2458891	30.872	ng/u1	0.00
49) Acenaphthylene-d8	14.387	160	2819110	31.310	ng/u1	0.00
54) 4-Nitrophenol-d4	14.893	143	360724	21.912	ng/u1	0.00
60) Fluorene-d10	15.681	176	2165926	32.292	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.804	200	258130	18.674	ng/u1	0.00
73) Anthracene-d10	17.557	188	3366225	32.719	ng/u1	0.00
81) Pyrene-d10	19.792	212	3604394	40.564	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.957	264	2364571	33.140	ng/u1	0.00

Target Compounds Qvalue

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

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