

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082823\
 Data File : BP016792.D
 Acq On : 28 Aug 2023 14:48
 Operator : MA/JU
 Sample : 04134-13
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCHK8

Quant Time: Aug 29 00:48:51 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	8.034	152	291270	20.000	ng/ul	0.00
20) Naphthalene-d8	10.863	136	1290011	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.687	164	823749	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.451	188	1806097	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	1373517	20.000	ng/ul	0.00
88) Perylene-d12	24.121	264	1215505	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	37770	5.359	ng/uL	0.00
4) Pyridine-d5	3.822	84	255703	11.873	ng/ul	0.00
7) Phenol-d5	7.169	99	170333	6.577	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.369	67	488049	29.566	ng/ul	0.00
11) 2-Chlorophenol-d4	7.552	132	457081	23.472	ng/ul	0.00
15) 4-Methylphenol-d8	8.734	113	309858	14.655	ng/ul	-0.01
21) Nitrobenzene-d5	9.228	128	291405	29.876	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.946	143	298723	29.055	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.469	165	506973	26.450	ng/ul	0.00
31) 4-Chloroaniline-d4	11.016	131	822730	27.990	ng/ul	-0.01
46) Dimethylphthalate-d6	14.098	166	1939222	31.216	ng/ul	0.00
49) Acenaphthylene-d8	14.387	160	2193781	31.239	ng/ul	0.00
54) 4-Nitrophenol-d4	14.892	143	45899	3.575	ng/ul	0.00
60) Fluorene-d10	15.681	176	1792414	34.263	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.798	200	239383	22.256	ng/ul	-0.01
73) Anthracene-d10	17.551	188	3114826	38.908	ng/ul	-0.01
81) Pyrene-d10	19.786	212	3455096	45.802	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.957	264	2390001	39.254	ng/ul	0.00

Target Compounds Qvalue

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

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