

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082624\
 Data File : BP021680.D
 Acq On : 27 Aug 2024 21:12
 Operator : RC/JU
 Sample : P3675-08
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCYE3

Quant Time: Aug 27 23:23:07 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP082324.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 27 05:11:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.798	152	302297	20.000	ng/u1	0.00
20) Naphthalene-d8	10.581	136	1250022	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.433	164	821034	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.233	188	1765363	20.000	ng/u1	-0.02
79) Chrysene-d12	21.674	240	1823248	20.000	ng/u1	-0.02
88) Perylene-d12	25.086	264	2100728	20.000	ng/u1	-0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.275	96	28947	3.179	ng/uL	0.00
4) Pyridine-d5	3.687	84	330664	12.555	ng/u1	0.00
7) Phenol-d5	6.957	99	555512	17.219	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	314632	17.921	ng/u1	0.00
11) 2-Chlorophenol-d4	7.328	132	366624	17.602	ng/u1	0.00
15) 4-Methylphenol-d8	8.492	113	436800	17.522	ng/u1	0.00
21) Nitrobenzene-d5	8.940	128	175251	17.426	ng/u1	0.00
24) 2-Nitrophenol-d4	9.663	143	176069	17.248	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	347620	17.280	ng/u1	0.00
31) 4-Chloroaniline-d4	10.710	131	519609	17.754	ng/u1	0.00
46) Dimethylphthalate-d6	13.839	166	1196444	19.045	ng/u1	0.00
49) Acenaphthylene-d8	14.122	160	1325550	18.793	ng/u1	0.00
54) 4-Nitrophenol-d4	14.633	143	190121	15.907	ng/u1	0.00
60) Fluorene-d10	15.445	176	1020034	19.313	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	112584	11.595	ng/u1	-0.01
73) Anthracene-d10	17.327	188	1635388	19.520	ng/u1	-0.02
81) Pyrene-d10	19.698	212	2021962	19.400	ng/u1	-0.02
92) Benzo(a)pyrene-d12	24.856	264	2227193	19.882	ng/u1	-0.02

Target Compounds	Qvalue

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

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