

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
 Data File : BP019066.D
 Acq On : 23 Dec 2023 07:21
 Operator : MA/JU
 Sample : 05869-01
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B26

Quant Time: Dec 24 23:57:49 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 21 04:46:12 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	204104	20.000	ng/u1	0.00
20) Naphthalene-d8	10.610	136	783880	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.451	164	513576	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.198	188	1202881	20.000	ng/u1	-0.01
79) Chrysene-d12	21.292	240	1245850	20.000	ng/u1	0.00
88) Perylene-d12	23.651	264	1469211	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	33546	5.601	ng/uL	0.00
4) Pyridine-d5	3.669	84	285270	17.781	ng/u1	0.00
7) Phenol-d5	6.993	99	632584	30.884	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	353070	29.084	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.346	132	491275	30.738	ng/u1	0.00
15) 4-Methylphenol-d8	8.534	113	513315	30.967	ng/u1	0.00
21) Nitrobenzene-d5	8.981	128	243771	34.940	ng/u1	0.00
24) 2-Nitrophenol-d4	9.699	143	280910	38.836	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	536001	36.239	ng/u1	-0.01
31) 4-Chloroaniline-d4	10.751	131	575077	29.751	ng/u1	-0.01
46) Dimethylphthalate-d6	13.869	166	1624543	35.442	ng/u1	0.00
49) Acenaphthylene-d8	14.145	160	1859992	36.851	ng/u1	0.00
54) 4-Nitrophenol-d4	14.669	143	279444	37.140	ng/u1	0.00
60) Fluorene-d10	15.445	176	1462863	38.021	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	262190	36.352	ng/u1	0.00
73) Anthracene-d10	17.298	188	2357624	37.282	ng/u1	0.00
81) Pyrene-d10	19.539	212	2940144	30.317	ng/u1	0.00
92) Benzo(a)pyrene-d12	23.498	264	3146882	36.735	ng/u1	0.00

Target Compounds	Qvalue

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

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