

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101821\
 Data File : BP007491.D
 Acq On : 19 Oct 2021 13:39
 Operator : CG/JU
 Sample : M4196-04
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JDGA3

Quant Time: Oct 19 14:13:56 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SFAM-EPA-BP101421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Oct 18 11:26:58 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.957	152	177586	20.00	ng/ul	0.00
20) Naphthalene-d8	10.763	136	725840	20.00	ng/ul	0.00
38) Acenaphthene-d10	14.592	164	470764	20.00	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	1013328	20.00	ng/ul	0.00
79) Chrysene-d12	21.433	240	1091076	20.00	ng/ul	0.00
88) Perylene-d12	23.880	264	1111154	20.00	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	19271	4.29	ng/uL	0.00
4) Pyridine-d5	3.805	84	42043	3.42	ng/ul	0.01
7) Phenol-d5	7.110	99	473117	34.11	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67	277048	33.49	ng/ul	0.00
11) 2-Chlorophenol-d4	7.487	132	374043	34.28	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	381198	35.68	ng/ul	0.00
21) Nitrobenzene-d5	9.110	128	178408	38.84	ng/ul	0.00
24) 2-Nitrophenol-d4	9.840	143	182466	42.53	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.381	165	367669	35.13	ng/ul	0.00
31) 4-Chloroaniline-d4	10.893	131	489485	34.59	ng/ul	0.00
46) Dimethylphthalate-d6	14.004	166	1169178	35.90	ng/ul	0.00
49) Acenaphthylene-d8	14.287	160	1428224	34.89	ng/ul	0.00
54) 4-Nitrophenol-d4	14.775	143	168275	32.20	ng/ul	0.00
60) Fluorene-d10	15.586	176	1004986	36.87	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	114373	28.17	ng/ul	0.00
73) Anthracene-d10	17.445	188	1603090	37.25	ng/ul	0.00
81) Pyrene-d10	19.674	212	1920716	35.21	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.721	264	1900071	34.37	ng/ul	-0.01

Target Compounds	Qvalue

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

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